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SEISMIC COUPLING FROM A NUCLEAR EXPLOSION
THE DEPENDENCE OF THE REDUCED DISPLACEMENT
POTENTIAL ON THE NONLINEAR REHAVIOR OF THE
NEAR SOURCE ROCK ENVIRONMENT

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SEISMIC COUPLING FROM A NUCLEAR EXPLOSION: THE DEPENDENCE OF THE REDUCED DISPLACEMENT POTENTIAL ON THE NONLINEAR BEHAVIOR OF THE NEAR SOURCE ROCK ENVIRONMENT

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Technical Report

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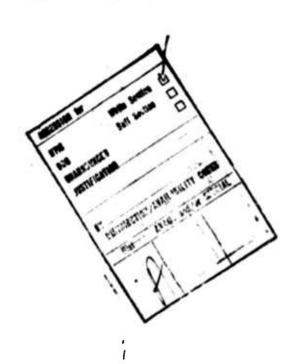
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A series of spherically symmetric calculations have been performed in which the properties of the near source material were varied systematically. Output from these calculations enables one to determine the relative teleseismic coupling efficiency of a given near source material. Also, since seismic magnitudes are directly related to explosive yield, this parameter study permits an assessment of uncertainties in the estimated yield in terms of uncertainties of the material properties of the test site.

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SUMMARY

During the past several years, Systems, Science and Software (S³) personnel have been actively engaged in a comprehensive program involving computer modeling of the nonlinear processes that characterize the near source ground motion from underground nuclear explosions, propagation of the resultant stress waves through the crust and upper mantle and prediction of the ground motion recorded at teleseismic distances from the explosion [Cherry, et al. (1973), Bache, et al. (1975)].

The objective of the research presented in this report is to determine the dependence of teleseismic magnitudes on the nonlinear behavior of the near source rock environment. Such information enables one to express uncertainties in explosive yield estimates in terms of uncertainties in the material properties and provides insight concerning the requirements for collection of geophysical data at a specific test site.

A series of spherically symmetric calculations have been performed in which the properties of the near source material were varied systematically. Output from these calculations enables one to determine the relative teleseismic coupling efficiency of a given near source material. Also, since seismic magnitudes are directly related to explosive yield, this parameter study permits an assessment of uncertainties in the estimated yield in terms of uncertainties of the material properties of the test site.

I. INTRODUCTION

1.1 OBJECTIVES

This report describes the results of a theoretical study directed at the determination of the sensitivity of the equivalent elastic source from a nuclear explosion to the nonlinear behavior of the near source rock environment. Uncertainties in yield estimates based on measured seismic magnitudes may be expressed in terms of uncertainties in the near source properties. The material parameters and the aspects of the constitutive model which have the greatest effect on the equivalent source are identified and guidelines are established for the collection of geophysical data in the near source region.

1.2 PROCEDURE

Equivalent elastic source calculations place severe requirements on both the constitutive model and the stress wave simulation code which contains the model. The constitutive model must account for the various modes of material behavior which occur between the cavity boundary and the distance at which the material response is elastic. The stress wave code must not only be able to easily accept the constitutive model but also accurately propagate the stress wave into the elastic region. The equivalent elastic source is then obtained from the properties of the displacement field in the elastic region.

Constitutive models, especially that portion of the model which pertains to the material strength, can be fairly complex. In order to isolate the effect of the constitutive model on seismic coupling we first assumed that only plastic flow would result when the allowable material strength was exceeded and an extensive parameter study was completed for

a water-tuif rock environment containing various amounts of air-filled porosity. We then modified the strength portion of the constitutive model to simulate tension failure, shear failure and the role of pore fluid pressure on effective stress.

Figure 1.1 depicts, in flow chart form, the organization of the parameter study. The material selected consists of a mixture of tuff and water which may contain air-filled voids. The properties of the water and grain density tuff constituents were held constant for all calculations. The mass fraction of water in the mixture, $f_{\rm w}$, is a natural parameter to select for initial variation and this choice is indicated in Fig. 1.1; the air-filled volume fraction, ϕ , was set to zero for this set of calculations. The equation of state for the tuff-water mixture is described in Section 2.1 of this report.

The remaining material parameters are varied one at a time while holding the water mass fraction constant at 0.17. Four of the material parameters are directly related to the porosity model which is discussed in Section 2.1: (1) the initial air-filled volume fraction, ϕ ; (2) the elastic limit of the pressure, $P_{\rm e}$; (3) the crush or compaction pressure, $P_{\rm c}$; (4) the bulk modulus, k. The shear modulus, μ , is selected as one of the independent elastic properties of the porous mixture. Given k, μ and the ambient density, ρ , the elastic P wave velocity, c, is given by

$$c = \sqrt{\frac{k_0 + \frac{4}{3} \mu}{\rho_0}}$$

while the shear wave velocity, β , is

$$\beta = \sqrt{\frac{\mu}{\rho_0}}$$

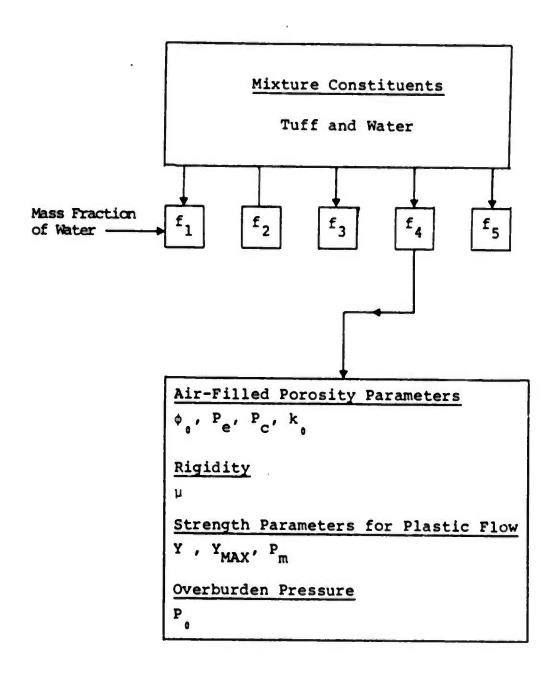


Figure 1.1. Flow chart for one-dimensional sensitivity study.

Additional parameters are contained in the material strength model which is discussed in Section 2.2. As noted earlier and shown in Fig. 1.1, only plastic flow was permitted during these calculations. The strength parameters included in this sensitivity study are: (1) the zero pressure yield strength, Y ; (2) the maximum yield strength, Y $_{\rm m}$; (3) the pressure at which the yield strength reaches its maximum value, P $_{\rm m}$.

One last parameter, the overburden pressure, P, is included in the one-dimensional sensitivity study although it depends on the depth of burial and is not a material property.

A detailed description of the material models employed in our near source calculations is presented in the next section. The results of the sensitivity study are discussed in Section III and conclusions based on those results are summarized in Section IV. In Section V the strength portion of the constitutive model is modified to include tension failure, shear failure and an effective stress law. The shape of the equivalent source spectrum is shown to be strongly dependent on both tension failure and the stress relaxation assumed to result from the effective stress formulation.

II. DESCRIPTION OF MATERIAL MODELS

For the parameter study we have chosen the rock constituent to be NTS tuff. The following sections, 2.1 and 2.2, present the pressure and material strength formulations used in the parameter study.

2.1 THE PRESSURE COMPONENT

Given the pressure component of the equation of state for both water and grain density tuff, then the pressure component for a fully saturated mix is obtained by specifying

$$f_W = \frac{M_W}{M_T + M_W} = \text{mass fraction cf water}$$
 (2.1)

Figures 2.1 and 2.2 show the Hugoniot and release isentropes for these two constituents.

At a given pressure \overline{P} a pressure equilibrium mix is obtained if

$$\overline{v} = \frac{v_r + f'v_w}{1 + f'} \tag{2.2}$$

$$e = \frac{e_r + f e_w}{1 + f'} \tag{2.3}$$

where

$$f' = \frac{f_w}{1 - f_w} = \frac{M_w}{M_r}$$
 (2.4)

 v_r , v_w , e_r and e_w are specific volumes and energies of the rock and water components of the pressure \overline{P} . It is now possible to generate a table that gives \overline{P} as a function of

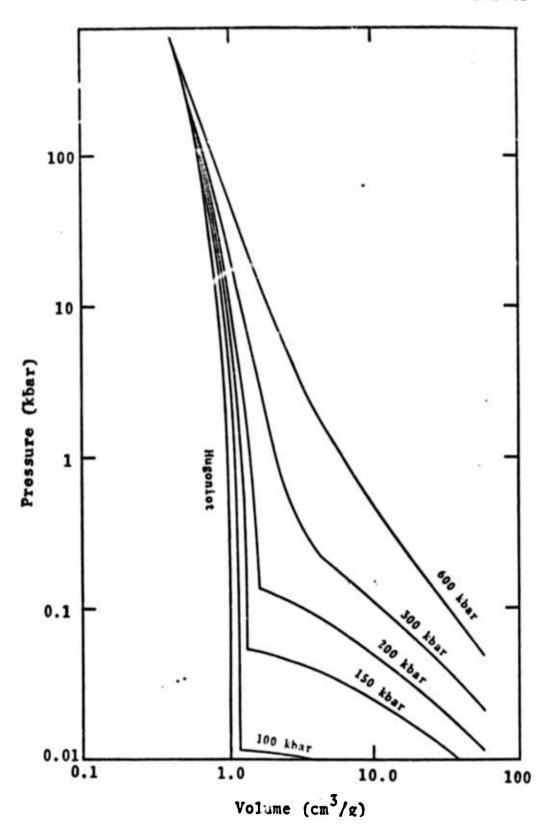


Figure 2.1. Hugoniot and release isentropes for water [Bjork and Gittings (1972)].

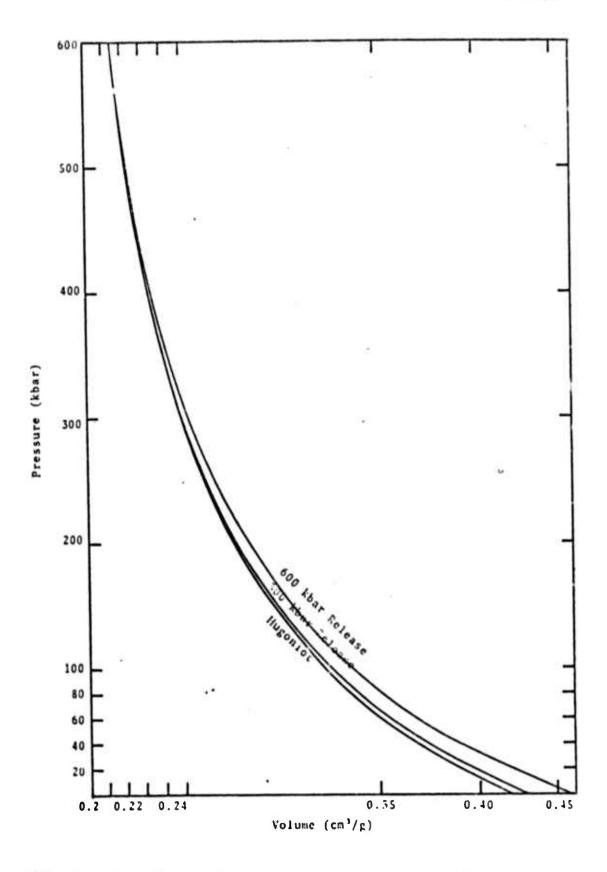


Figure 2.2. Hugoniot and release isentropes for tuff, ρ = 2.4 g/cc [Riney, et al. (1972)].

 \overline{v} and e. This table (Riney, et al.) [2] suitably chooses values of \overline{v} and e so that no time consuming search is needed when performing an interpolation at a given thermodynamic state. Figure 2.3 shows the result of mixing the rock and water constituents shown in Figs. 2.1 and 2.2 using a 0.17 value for f_w .

Air-filled porosity is included in the model by defining a parameter α , given by

$$\alpha = \frac{\mathbf{v_r} + \mathbf{v_w} + \mathbf{v_p}}{\mathbf{v_r} + \mathbf{v_w}} = \frac{\mathbf{v}}{\overline{\mathbf{v}}}$$
 (2.5)

where v is the specific volume of the partially saturated mix. Then

$$P = \frac{1}{\alpha} \overline{P} (v/\alpha, e)$$
 (2.6)

where \overline{P} is found from the table using v, e. Equation (2.6) is equivalent to the porosity model proposed by Herrmann (3) except \overline{P} is weighted with $1/\alpha$. Both v and e are zone variables and are obtained from SKIPPER at any given time cycle.

The parameter α is assumed to vary in the following way during loading (pore collapse) [Cherry, et al.] [4]

$$\alpha = 1 \quad P > P_{C} \tag{2.7a}$$

$$\alpha = 1 + (\alpha_e^{-1}) \left(1 - \frac{P - P_e}{P_c^{-P_e}}\right) \quad P_e \leq P \leq P_c \quad (2.7b)$$

$$\alpha = \alpha_0 + \beta \left(1 - e^{n P/P} e\right) 0 \le P \le P_e$$
 (2.7c)

$$\alpha = P\alpha'_{p=0} + \alpha P < 0 \quad (\alpha' = d\alpha/dp)$$
 (2.7d)



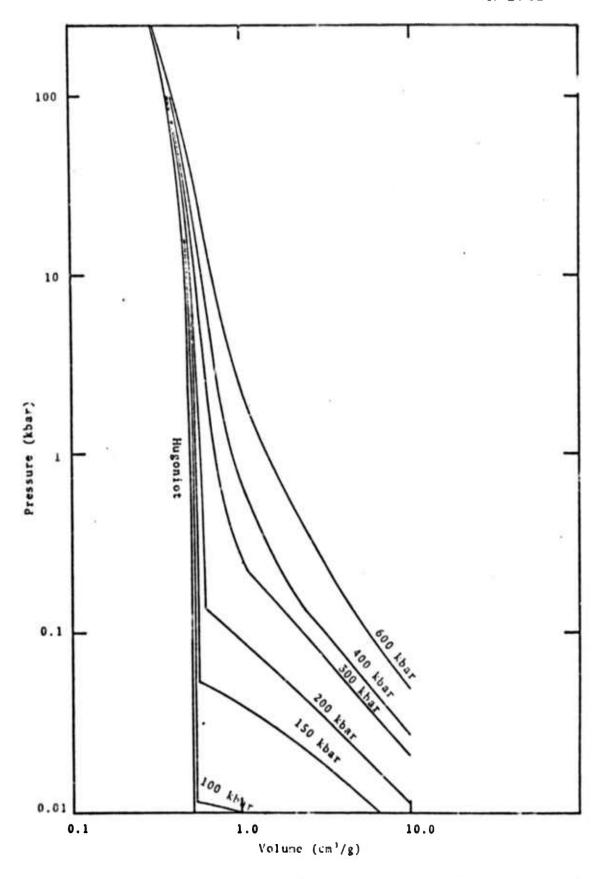


Figure 2.3. Hugoniot and release isentropes for saturated tuff (f = 0.17).

where P_e and P_c are the elastic and crush pressure. β and η are determined by specifying the zero pressure bulk modulus (k₀) of the porous mix and requiring that α be continuous at $P = P_e$. The initial air-filled volume fraction (ϕ ₀) and α ₁ are related by

$$\phi_0 = \frac{\alpha - 1}{\alpha_0} = \frac{V_p}{V_w + V_r + V_p}$$
 (2.8)

while the volume fractions of rock and water $(\phi_{\mathbf{r}} \text{ and } \phi_{\mathbf{w}})$ are given by

$$\phi_{\mathbf{r}} = \frac{1 - \phi}{1 + \frac{1}{A}} \tag{2.9}$$

$$\phi_{W} = \left(\frac{1 - \phi}{1 + A}\right) A \tag{2.10}$$

where

$$A = \frac{\rho_r}{\rho_w} \frac{f}{1-f}$$

$$\rho_r = 2.4 \text{ g/cc}$$

$$\rho_w = 1.0 \text{ g/cc}$$

We have found (Cherry, et al. [4]) that a pore recovery formulation needed to be included in the model in order to insure a smooth transition between loading states near P_e . If α^* is the minimum α seen by a zone then

$$\alpha = t\alpha^* + (1-t) \alpha (P)$$
 (2.11)

where (P) is given by Eqs. (2.8) and

$$t = 0, \alpha^* \ge \alpha_e = \frac{\alpha^* - \alpha_e}{1 - \alpha_e}, 1 \le \alpha^* \le e$$
 (2.12)

This assures full recovery when $\alpha^* = \alpha_e$ and zero recovery when $\alpha^* = 1$.

Figure 2.4 shows the loading and release P-v curves that result from adding 5 percent air-filled porosity to the mix of Fig. 2.3. We assume $P_e = 0.15$ kbar and $P_c = 1.25$ kbar. Also shown for comparison in the figure are the hydrostatic unloading data for unit 3 tuff at the Diamond Dust site (Stephens, et al. [5]).

In summary, the pressure component of the equation of state is uniquely determined by assuming a pressure equilibrium mix and specifying f (Eq. (2.1)), ϕ (Eq. (2.8)), P_e , F_c , k (Eq. (2.7b) and (2.7c)) and the grain density of the rock component.

2.2 MATERIAL STRENGTH

A strength relation of the form

$$S_{ij} S_{ij} \le \frac{2}{3} Y^2 (P_T, e)$$
 (2.13)

where S_{ij} is the stress deviator, $P_T = P + P_0$, P_0 equals the overburden pressure, and

$$Y(P,e) = \left[Y_0 + Y_m \frac{P}{P_m} (2 - \frac{P}{P_m})\right] (1 - \frac{e}{e_m}) P < P_m$$

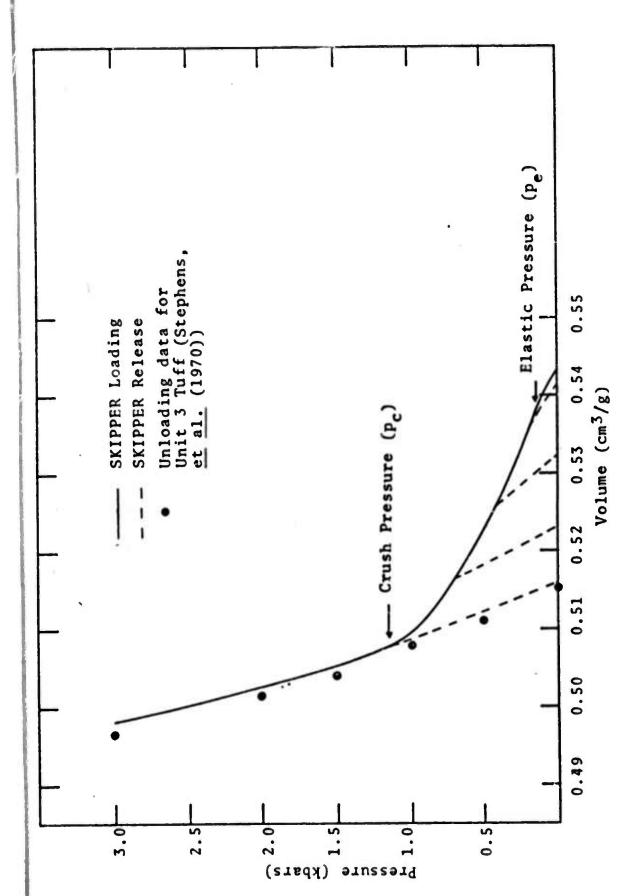
$$= (Y_0 + Y_m) (1 - \frac{e}{e_m}), P \ge P_m \qquad (2.14)$$

$$= 0 \qquad e \ge e_m$$

For spherical symmetry Eq. (2.13) becomes

$$|S_1| \le \frac{2}{3} Y (P_T, e)$$
 (2.15)

For all the calculations performed for the parameter study, a generalized associated flow rule was used for the material in which the volumetric plastic strain rate was



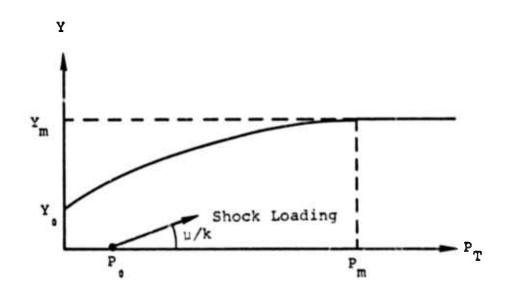
Loading and release p-V curves for partially saturated tuif (f = 0.17, $\phi_0 = 0.05$) [Cherry, et al. (1973)]. Figure 2.4.

assumed to be zero. This assumption implies (Cameron and Scorgie^[6]) that if $|S_1|$ exceeds $\frac{2}{3}$ Y, then

$$S_{1} = \left(\frac{2}{3} Y\right) \text{ sign } (S_{1})$$
 (2.16)

The material strength dependence on P_0 , Y_0 , Y_m and P_m is summarized in Fig. 2.5. The shock loading path is also shown in the figure to have a slope of μ/k where k is the effective bulk modulus of the shock wave. This loading path is only valid for regions where the radial strain is much greater than the tangential strain, i.e., at the shock front.

In Section V the material strength formulation will be modified to include brittle failure and an effective stress law.



$$Y = \left(Y_0 + Y_m \frac{P_T}{P_m}\right) \left(2 - \frac{P_T}{P_m}\right) \qquad P_T < P_m$$

$$Y = Y_0 + Y_m \qquad P_T > P_m$$

Y_n = Maximum allowable stress difference

P_T = Hydrodynamic component of the stress tensor including overburden pressure

P = Overburden pressure

Figure 2.5. Assumed relationship between the material strength (Y) and the hydrodynamic component of stress (P_T).

III. DISCUSSION OF RESULTS

3.1 THEORETICAL CONSIDERATIONS

A detailed discussion of the reduced displacement potential, RDP, which serves as the equivalent elastic source representation for spherically symmetric explosions is presented in Appendix A. The SKIPPER code, which is used to propagate the shock wave into the elastic region is discussed in Appendix B. The results presented in Fig. A.1 show the calculated reduced velocity potential spectra using the material model described in the preceding section. As noted in Appendix A, the spectra are essentially flat in the teleseismic frequency band for device yields at least as large as 200 kt. This result along with Eq. (A.9) imply that body wave magnitude, m_b, has the proportionality

$$m_{b} \sim \log \left[c\Psi(\infty)\right]$$
 (3.1)

where $\Psi(\infty)$ is the steady state value of the RDP and c is the near source compressional wave speed. Of course, this relation does not include the effect of pP or any of the details of the geophysical properties of the crustal layers. It has been derived by considering the impedence mixmatch between the low velocity source region and the basement rock, i.e., a two layer earth model. It is prefered to the homogeneous earth assumption which gives $\Psi(\infty)/\mathbb{C}$ for m_b proportionality.

If $\Psi_i^{(\infty)}$ and $\Psi_k^{(\infty)}$ are two values of $\Psi^{(\infty)}$ corresponding to materials i and k then the change in teleseismic magnitude is g_ven by

$$\Delta m = m^{i} - m^{k} = \log \left[\frac{c_{i} \Psi_{i}(\infty)}{c_{k} \Psi_{k}(\infty)} \right]$$
 (3.2)

If material properties remain constant and the only change is the device yield, W, then cube root scaling gives

$$\Delta m = m^{i} - m^{k} = \log \frac{W_{i}}{W_{k}}$$
 (3.3)

Equation (3.2) gives the "scaling" law for teleseismic magnitudes in terms of $c\Psi(\infty)$. This equation is always valid as long as the spectrum of the equivalent source is flat within the teleseismic frequency band and path effects associated with earth structure are invariant between events.

It should be noted that uncertainties in yield estimates based on experimentally determined seismic magnitudes may be related to uncertainties in near source material properties via direct application of Eq. (3.2).

3.2 PARAMETER SENSITIVITY

The results of the sensitivity study are summarized in Table 3.1. Some general remarks are in order before discussing the effect of individual parameter variations. Note that the ambient density, ρ , and the bulk modulus, k , are included in the table although they are not independent parameters in the context of this study; i.e., their values are determined by the independent parameters appearing in the first ten columns. Calculation No. 18 is selected as the reference for body and surface wave magnitude variations in order to avoid the appearance of minus signs in the Δm_h column. Calculations marked with an asterisk in the $\Psi(\infty)$ column were performed with a new source description which gave an RDP 42 percent higher than the RDP obtained in Calculation No. 8 using the old source description; the RDP's calculated with the new source description were therefore scaled down by a factor of 0.704 and the scaled values are reported in Table 3.1. All calculations appearing in Table 3.1 were performed with a device yield of 20 tons.

TABLE 3.1
RESULTS OF PARAMETER SENSITIVITY STUDY

ي	••	P. (Johan)	P. (thar)	C (Dar/sec)	u (thar)	Y. (Other)	Y, MAX (Abar)	P m ()dar)	P. (kbar)	(gm/cm³)	k (kbar)	f.	€ °	Calculation
0			1	3.81	40.0	0	0.5	9.0	0.116	2.4	294.4	13.88		1
0.05				2.80	40.0	0	0.5	9.0	0.116	2.243	121.8	10.01		7
0.136				2.44	40.0	0	0.5	9.0	0.116	2.015	66.5	7.85	_	9
0.17				2.40	40.0	0	0.5	9.0	0.116	1.938	57.8	7.34	-	•
0.19				2.38	40.0	c	0.5	9.0	0.116	1.895	53.8	7.09	_	\$
0.17				2.40	40.0	0	0.5	9.0	0.116	1.90	56.1	3.60	_	٠
0.17				2.40	40.0	0	0.5	9.0	0.116	1.84	52.6	2.10	_	7
0.17				2.40	40.0	0	1.0	9.0	0.116	1.84	9::5	1.35	-	∞
0.17				2.40	40.0	0	1.0	9.0	0.116	1.84	52.6	1.48*	_	'n
0.17				2.40	40.0	0	0.5	9.0	0.075	1.90	\$5.1	4.20		10
0.17			1.25	2.40	40.0	•	0.5	9.0	0.075	1.90	56.1	4.60	0.6059	11
0.17				2.32	10.0	0	1.0	9.0	0.116	1.84	45.7	1.31		12
0.17				2.25	40.0	0	1.0	9.0	0.116	1.84	39.8	1.29	_	13
0.17				2.40	42.0	0	1.0	9.0	0.116	1.84	50.0	1.28*		14
0.17				2.40	45.0	0	1.0	9.0	0.116	1.84	46.0	1.17		15
0.17				2.40	40.0	0.1	1.0	9.0	0.116	1.84	52.6	1.32		16
0.17				2.40	40.0	0.5	1.0	9.0	0.116	1.84	52.6	1.31		11
0.17				2.40	40.0	1.0	1.0	9.0	0.116	1.84	52.6	1.14		18
0.17				2.40	40.0	0	9.5	1.25	0.075	1.90	56.1	7.70		19
0.17				2.40	40.0	0	0.5	9.0	0.348	1.90	56.1	2.41		20
0.17				2.40	40.0	•	9.9	9.0	0.619	1.90	56.1	2.08		77
0.17	9.02	0.075			40.0	•	0.75	9.0	0.116	1.84	52.6	1.51*		8

Calculation performed with new source description.

The effects of water content on seismic coupling are revealed by Calculations 1 through 5. The reduced displacement potential was calculated for five fully saturated tuffs having different mass fractions of water, f_{ω} . Fully saturated tuffs were chosen in order to eliminate the effects of airfilled porosity. Note that adding water decreases both the density, ρ , and the bulk modulus of the mixture, k. Since the shear modulus is held constant at 40 kbars, variations in compressional wave speed, c, reflect only the changes in density and bulk modulus. Variations in body wave magnitude as functions of water mass fraction are described in Fig. 3.1. The decrease in seismic coupling as the water content is increased is to be expected. Since water is more compressible than rock at all pressures of interest, the amount of energy dissipated in the form of heat via shock loading increases with increasing water content, leaving less energy available for conversion into seismic radiation. mb decreases by about 0.5 magnitude units as the mass fraction of water increases from 0 to 0.19.

The effect of air-filled voids on seismic magnitude is obtained from Calculations 4, 6 and 7. The water content is held constant at $f_w = 0.17$ for these and the remaining calculations in Table 3.1. The variation of seismic magnitude as a function of the void fraction, ϕ_0 , is presented in Fig. 3.2. Note that the introduction of air-filled voids causes a drastic reduction in seismic coupling and the importance of establishing this parameter preshot is obvious.

Calculations 8 and 9 show that seismic magnitude is not sensitive to P_e , the pressure below which pore closure is elastic and fully recoverable. A factor of two increase in P_e increases the seismic magnitude by 0.04 units. Similarly, Calculations 10 and 11 show that magnitude is insensitive to variations in P_c , the pressure at which all ambient porosity is irreversible removed. An increase in P_c by a factor of

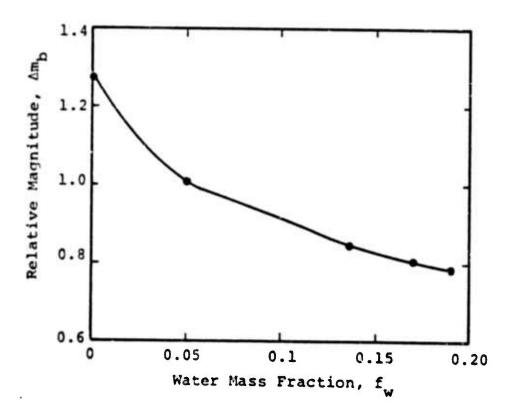


Figure 3.1. Variation of seismic magnitudes with water mass fraction.

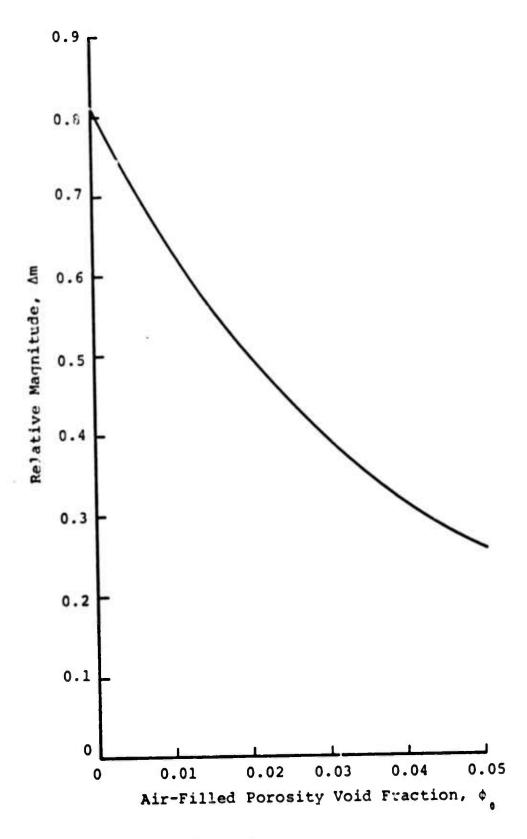


Figure 3.2. Effect of air-filled porosity on seismic magnitude.

2.5 produces an increase in magn_tude of 0.03 units. It should be noted that P_e and P_c are associated with the porosity model and might increase in importance as ϕ , the air-filled void fraction, is increased.

The effect of variations in near source seismic velocity, c, on magnitude is obtained from Calculations 8, 12 and 13. The permissible range of c is limited at this point because the properties of tuff, the properties of water, the mass fraction of water and the shear modulus are held constant. Variations in c therefore reflect only a change in the initial slope of the crush curve associated with the porosity model. As shown in F² 3, a 6 percent error in P wave velocity produces a cha. m_b of about 0.05 magnitude units. Increasing the compressional wave speed increases seismic coupling for body waves.

Calculations 8, 14 and 15 illustrate the effect of variations in the shear modulus on seismic magnitude. Body wave magnitude is reduced as the shear modulus increases as shown in Fig. 3.4.

The yield surface is described by the parameters Y, Y_{m} , and P_{m} . Calculations 8, 16, 17 and 18 show the effect of $Y_{\underline{a}}$, the zero pressure yield strength, on seismic coupling. The variation of Δm with Y is plotted in Fig. 3.5. Increases in Y tend to reduce seismic coupling. The results of Calculations 7, 8 and 22 are plotted ir. Fig. 3.6 which shows the effect of $Y_{\mbox{\scriptsize MAX}}$, the maximum allowable material strength, on seismic magnitude. Increasing Y_{MAX} by a factor of two causes magnitude to decrease by 0.192 magnitude units, indicating that $Y_{\mbox{\scriptsize MAX}}$ is a fairly sensitive parameter. For a given value of Y_{MAX} , the shape of the material strength surface is controlled by P_{m} , the pressure at which Y_{MAX} is attained. From Calculations 10 and 19 we see that increasing $P_{\rm m}$ by a factor of 2.98 causes magnitude to increase by 0.26 units. This result indicates that the shape of the failure surface plays an important role in determining

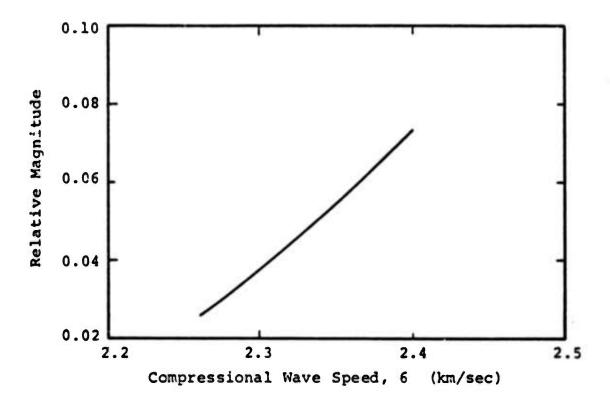


Figure 3.3. Effect of compressional wave speed on seismic magnitude.

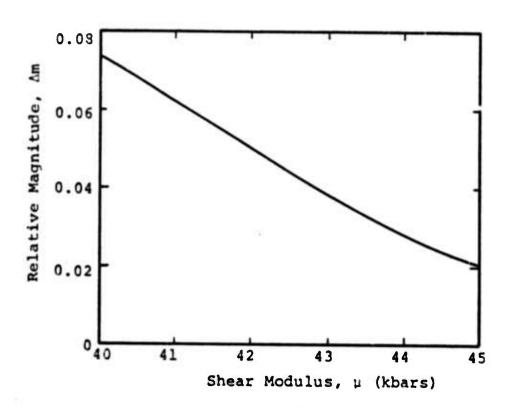


Figure 3.4. Effect of shear modulus on seismic magnitude.

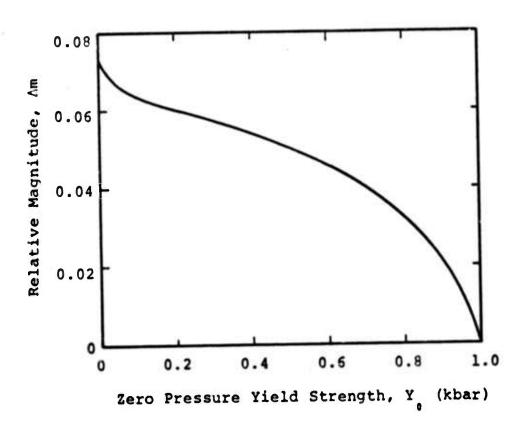


Figure 3.5. Effect of zero pressure yield strength on seismic coupling.

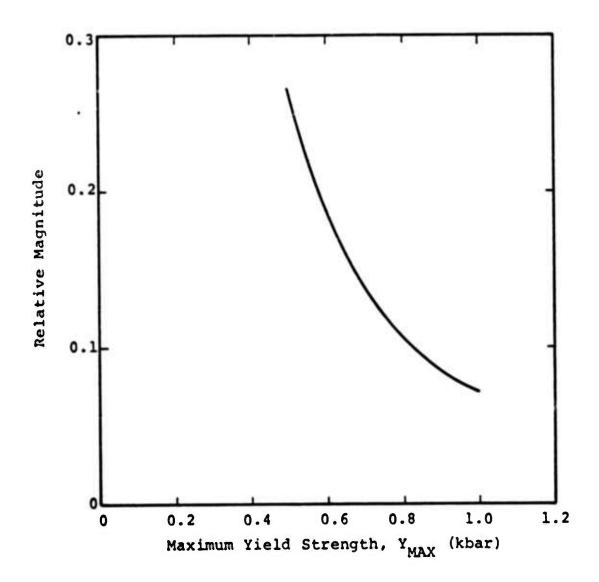


Figure 3.6. Effect of maximum material strength on seismic magnitude.

coupling efficiency. The material has less strength at low pressures when P_m is increased and we should expect a corresponding increase in magnitude. In fact, if any of the parameters Y_0 , Y_m , or P_m are varied such that the material strength is enhanced, a decrease in coupling efficiency is observed.

The effect of overburden pressure, P , on magnitude, as determined by Calculations 6, 10, 20 and 21, is shown in Fig. 3.7. If all material properties are held constant, coupling efficiency decreases with increasing overburden pressure. An increase in the depth of burial, DOB, by a factor of six causes magnitude to decrease by a factor of two. The decrease in magnitude is not uniform with depth. An increase in DOB at shallow depths causes magnitude to decrease more than that caused by an equivalent increase in DOB at greater depths. This effect results from the rapid increase in material strength at low pressures. We should note that it is unlikely that material properties would remain invariant as overburden pressure is increased from 100 bars to 600 bars. An increased DOB usually corresponds to an increase in water content, resulting in decreased airfilled porosity and material strength.

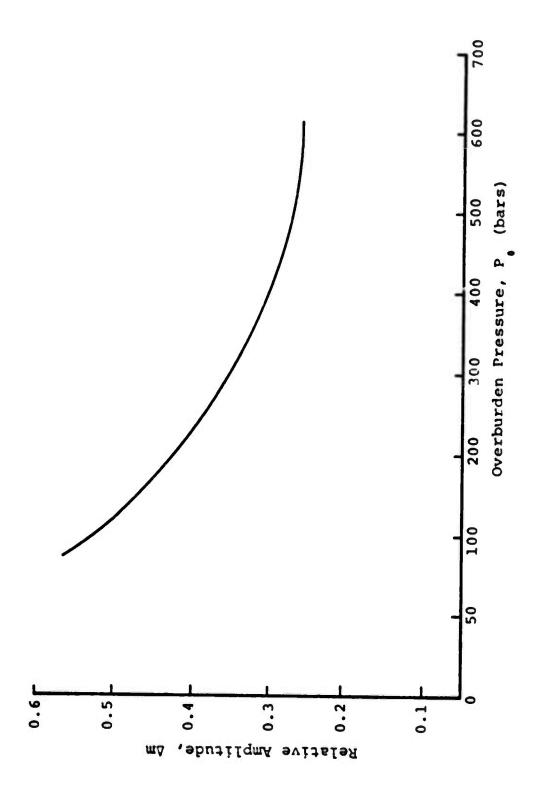


Figure 3.7. Effect of overburden pressure on seismic magnitude.

IV. SUMMARY OF RESULTS

The results of the parameter study provide valuable insight concerning the importance of various material parameters with respect to accurate prediction of seismic coupling. has been shown that increasing the overburden pressure, or, equivalently, the depth of burial, decreases coupling efficiency substantially. Therefore, the DOB must be accurately known in order to make a reliable prediction of seismic magnitude. It is also desirable to know the properties of the material in the immediate vicinity of the device; if substantial spatial variation of the material properties occurs near the device or if the DOB is quite shallow it may be necessary to consider two-dimensional effects which are not treated here. Accurate prediction of seismic magnitudes via one-dimensional spherically symmetric calculations is therefore understood to be dependent on the material uniformity about the device.

Since the elastic properties of the near source material appear in the magnitude relation, Eq. (3.2), via the P wave velocity, c, it is obviously important to determine their values accurately. A positive error in determination of c would cause body wave magnitude to be over-predicted. A positive error in μ would cause magnitude to be underpredicted.

The steady state value of the RDP, $\Psi(\infty)$, which appears in the magnitude relation, Eq. (3.2), is dependent on the shock response of the near source material. The near source material properties which have the most pronounced effect on the shock response are the water mass fraction, f_{ψ} , and the air-filled void fraction, ϕ . Positive errors in either f_{ψ} or ϕ could lead to substantial under-prediction of seismic magnitudes. Seismic magnitude is not very sensitive

to $P_{\rm e}$, the elastic pressure, and $P_{\rm c}$, the crush pressure, indicating that details in the porosity model are relatively unimportant.

Seismic magnitudes are, however, very sensitive to the parameters which describe the failure surface. If $Y_{\mathfrak{g}}$, $Y_{\mathfrak{m}}$, or $P_{\mathfrak{m}}$ are varied such that the material strength is enhanced the coupling efficiency of an explosive device is impaired. Thus, a positive error in the material strength would lead one to an under-prediction seismic magnitude.

V. THE EFFECT OF MATERIAL STRENGTH ON THE PRESSURE AND DEVIATORIC COMPONENTS OF THE STRESS TENSOR

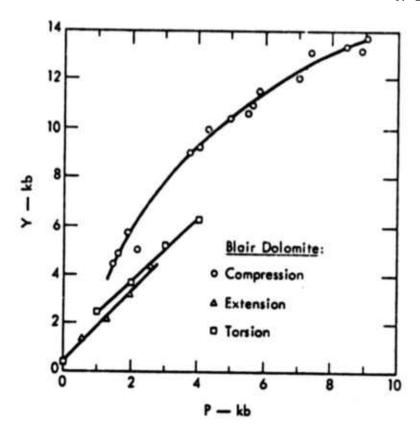
5.1 INTRODUCTION

Most code calculations of the type presented in Section 3, either for ground shock predictions or equivalent source calculations, have assumed that material strength depends only on the pressure component of the stress tensor (i.e., Eq. (2.14)). Also, very little attention has been paid to the consequences of exceeding the allowable material strength other than envoking either an associated or non-associated flow rule, i.e., Eq. (2.16). Further, the influence of water content on material strength via an effective stress law has not been considered in any ground shock calculation performed to date.

A number of investigators have considered the above topics separately. In particular, Cherry and Petersen [7] have identified the dependence of material strength on the third deviatoric invariant. Maenchen and Sack [10] presented a formulation for tension failure which is ideally suited for a stress wave code. Garg and Nur [11] investigated the consequences of various effective stress laws on the interpretation of experimental data. We have relied heavily on the results of these investigators in order to formulate the strength portion of the constitutive model.

5.2 THE DEPENDENCE OF MATERIAL STRENGTH ON THE STRESS IN-VARIANTS

Figure 5.1 is taken from Cherry and Petersen^[7] and shows the improvement in the definition of material strength, Y, when P is replaced by \overline{P} . Similar improvement has also been found for granite, limestone and glass. Y, P and \overline{P} are obtained from the stress invariants J_1 , J_2 , J_3 , as follows:



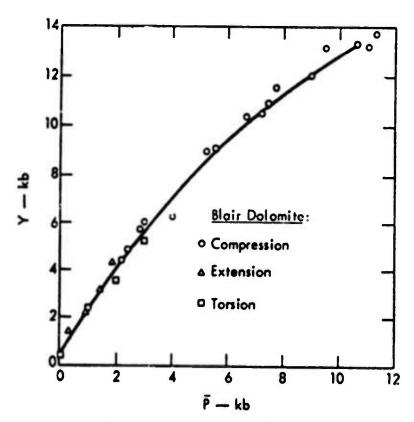


Figure 5.1. These figures are from Cherry and Petersen [1970] and show the improvement in failure surface definition when P is replaced by P. The strength data are from Handin, et al. [1967].

$$Y = (3J_2)^{1/2} (5.1)$$

$$\overline{P} = P - \frac{1}{2} \left(\frac{J'}{2} \right)^{1/2} \tag{5.2}$$

$$P = -\frac{J}{3} \tag{5.3}$$

P is now the pressure component of the stress tensor discussed in Section II with the overburden pressure, P, added. J and J are the first, second deviatoric and third deviatoric stress invariants. If σ_1 , σ_2 , σ_3 are principal stresses, then

$$J = \sigma + \sigma + \sigma$$
1 1 2 2 3 3 (5.4)

$$J_{2}^{\prime} = \frac{(\sigma_{11} + P)^{2} + (\sigma_{12} + P)^{2} + (\sigma_{13} + P)^{2}}{2}$$
 (5.5)

$$J' = (\sigma_1 + P) (\sigma_2 + P) (\sigma_3 + P)$$
 (5.6)

Note that when the intermediate principal stress, σ_{22} , is equal to either the maximum, σ_{11} , or minimum, σ_{33} , principal stress, then

$$Y = \left| \sigma_{11} - \sigma_{22} \right| \tag{5.7}$$

$$\overline{P} = -\left(\frac{\sigma_1 + \sigma_2}{2}\right) \tag{5.8}$$

The functional form we have chosen to represent the variation of material strength, Y, with the stress state, \overline{P} , is identical to that used in Section 2, i.e.,

$$Y = \left\{ Y_0 + \frac{Y_m}{\overline{P_m}} \quad \overline{P} \left[2 - \frac{\overline{P}}{\overline{P_m}} \right] \right\} \left[1 - \frac{e}{e_m} \right] \quad \overline{P} < P_m$$

$$Y = (Y_0 + Y_m) \left[1 - \frac{e}{e_m} \right] \quad \overline{P} > \overline{P_m}$$

$$Y = 0 \quad e > e_m$$

$$(5.9)$$

where Y, Ym, \overline{P}_m and e_m are input constants. These constants along with P, the overburden pressure, determine the variation of material strength with stress state along with the initial stress state, P, corresponding to the overburden pressure at the depth of burial of the device.

The material strength dependence on P , Y , Y and \overline{P}_m is summarized in Fig. 5.2. The shock loading path is also shown in the figure to have a slope of 1-2 σ , where σ is Poisson's ratio. This loading path is only valid for regions where the radial strain is much greater than the tangential strain, i.e., at the shock front.

5.3 SHEAR FAILURE

Hooke's law is used to obtain an initial estimate of the stress deviators, i.e.,

$$\hat{s}_{ij}^{n+1} = s_{ij}^{n} + 2\mu \, \dot{e}_{ij}^{n} \, \Delta t$$
 (5.10)

where \hat{S}_{ij}^{n+1} and S_{ij}^{n} are the values of the stress deviator at time t + Δt and t respectively.

μ is the shear modulus

eij is the strain rate devaitor

At is the time increment

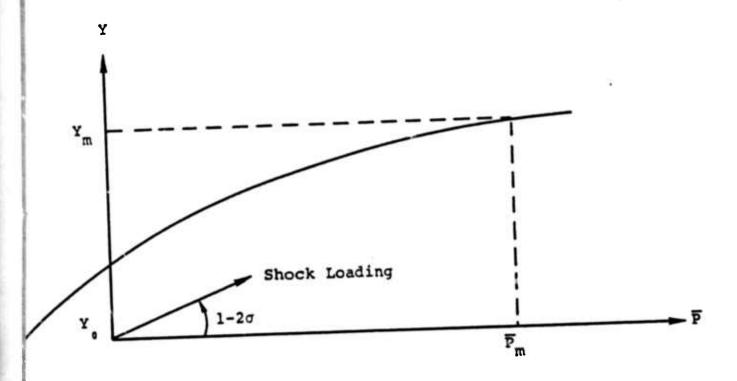


Figure 5.2. Dependence of Y on Y , \overline{P}_m and Y_m . When \overline{P} is used for the abcissa. The shock loading path has a slope of 1-2 σ .

Shear failure occurs if the material strength, given by Eq. (5.9) and evaluated at \hat{P} , is exceeded, i.e., if

$$J > Y(\widehat{P}) \tag{5.11}$$

where

$$J = \left[\frac{3}{2} \left(\hat{S}_{11}^2 + \hat{S}_{22}^2 + \hat{S}_{33}^3\right)\right]^{1/2}$$
 (5.12)

$$\hat{\bar{p}} = p^{n+1} - \frac{1}{2} \left(\frac{\hat{s}_{11} \hat{s}_{22} \hat{s}_{33}}{2} \right)^{1/3}$$
 (5.13)

 \mathbf{p}^{n+1} includes the overburden pressure and we have omitted the n+1 superscript on $\hat{\mathbf{S}}_{i,i}$.

If $J > Y(\overline{P})$ then adjustment of the stress deviators, \hat{s}_{ij}^{n+1} , is required. We assume that

$$s_{ij}^{n+1} = a \hat{s}_{ij}^{n+1}$$
 (5.14)

where S_{ij}^{n+1} is the adjusted value of the stress deviator and

$$a = \frac{Y(\hat{\overline{P}}) + \frac{b}{2} \left(\frac{\hat{S}_{1} + \hat{S}_{2} + \hat{S}_{3}}{2} \right)^{1/3}}{J + \frac{b}{2} \left(\frac{\hat{S}_{1} + \hat{S}_{2} + \hat{S}_{3}}{2} \right)^{1/3}}$$
(5.15)

$$b = \frac{dY}{d\overline{P}} \quad \text{(evaluated at } \hat{\overline{P}}\text{)} \tag{5.16}$$

Equation (5.15) is obtained by approximating the strength function at \hat{P} with a first order Taylor series and assuming that no adjustment in P^{n+1} occurs during shear failure.

5.4 TENSION FAILURE

Tension failure is assumed to occur in the element if a principal stress is greater than zero and if the inequality of (5.11) has ever been satisfied. We then apply the tension failure model proposed by Maenchen and Sack [10] and introduce an inelastic strain normal to the crack. This inelastic strain is just sufficient to zero the tensile stress.

For example, if $\hat{\sigma}_{11}$, $\hat{\sigma}_{22}$ and $\hat{\sigma}_{33}$ are the three principal stresses and if $\hat{\sigma}_{11}$ is greater than zero, then the adjusted stress (σ_{11} , σ_{22} , σ_{33}) are given by

$$\sigma_{11} = \hat{\sigma}_{11} - (k + \frac{4}{3} \mu) \Delta e_{11}$$

$$\sigma_{22} = \hat{\sigma}_{22} - (k - \frac{2}{3} \mu) \Delta e_{11}$$

$$\sigma_{33} = \hat{\sigma}_{33} - (k - \frac{2}{3} \mu) \Delta e_{11}$$
(5.17)

where

$$\Delta e_{11} = \frac{\hat{\sigma}}{k + \frac{4}{3} \mu}$$
 (5.18)

k is the bulk modulus, $\;\mu\;$ is the shear modulus and all stresses include the overburden pressure.

If two of the principal stresses, say σ and σ are greater than zero, then the stress adjustment becomes

$$\sigma_{11} = \hat{\sigma}_{11} - (k - \frac{2}{3} \mu) (\Delta e_{11} + \Delta e_{22}) - 2\mu \Delta e_{11}$$

$$\sigma_{22} = \hat{\sigma}_{22} - (k - \frac{2}{3} \mu) (\Delta e_{11} + \Delta e_{22}) - 2\mu \Delta e_{22}$$

$$\sigma_{33} = \hat{\sigma}_{33} - (k - \frac{2}{3} \mu) (\Delta e_{11} + \Delta e_{22})$$

$$(5.19)$$

where

$$\Delta e_{11} = \frac{(k + \frac{4}{3}) \hat{\sigma}_{11} - (k - \frac{2}{3}\mu) \hat{\sigma}_{22}}{4\mu (k + \frac{\mu}{3})}$$

$$\Delta e_{22} = \frac{(k + \frac{4}{3}\mu) \hat{\sigma}_{22} - (k - \frac{2}{3}\mu) \hat{\sigma}_{11}}{4\mu (k + \frac{\mu}{3})}$$
(5.20)

All the inelastic strain increments (Δe_{11} , Δe_{22} , Δe_{33}) are accumulated on each cycle giving

$$E_{11}^{n+1} = E_{11}^{n} + \Delta e_{11}$$

$$E_{22}^{n+1} = E_{22}^{n} + \Delta e_{22}$$

$$E_{33}^{n+1} = E_{33}^{n} + \Delta e_{33}$$
(5.21)

These equations give the basic stress-strain adjustment during tension failure. However, they apply equally well for crack closure. If at least one of the total strains in Eq. (5.21) (say \mathbf{E}_{11}^{n}) is greater than zero, then the crack will open or close depending on the sign of $\hat{\sigma}_{11}$. If $\hat{\sigma}_{11} > 0$, then

$$\Delta e_{11} > 0$$

$$E_{11}^{n+1} > E_{11}^{n}$$

and the crack width increases. The inequalities are reserved if \hat{T} < 0 and the crack width decreases. Closure will continue until

$$E_{11}^{n} + \Delta e_{11} < 0$$
.

Then

$$\Delta e_{11} = - E_{11}^{n}$$

$$E_{11}^{n+1} = 0$$

and the crack is completely closed. When this state is achieved the element is able to support a compressive stress in the (1,0,0) direction.

5.5 EFFECTIVE STRESS

All effective stress laws, formulated for fluid-saturated porous rocks, assume that the fluid pressure should be included in the calculation of the stress state which determines material strength. An excellent review of various definitions of effective stress and their consequences in terms of laboratory strength data has been given by Garg and Nur. [11]

The definition of stress state we have adopted is \overline{P} , where \overline{P} is given by Eq. (5.2). If P_p is the pore fluid pressure, then the effective stress state becomes

$$\overline{P} = P - \frac{1}{2} \left(\frac{J'}{2} \right)^{1/3} - F(P_p)$$

where $F(P_p)$ is a positive function of the pore fluid pressure. We have assumed that $F(P_p)$ behaves as follows:

$$F(P_p) = 0$$
 if tension cracks are open. (5.22)

$$F(P_p) = 0$$
 if P has never exceeded the crush pressure, P_c . (5.23)

$$F(P_p) = 0 \text{ if } P - \frac{1}{2} \left(\frac{J}{2}\right)^{1/3} < 0.$$
 (5.24)

If none of the above apply, then

$$F(P_p) = P - \frac{1}{2} \left(\frac{J'}{2}\right)^{1/3}$$
 (5.25)

The conditions given in Eqs. (5.22) and (5.23) assume that the pore fluid pressure has no effect on the stress state when air filled porosity exists in the element. If no air filled porosity exists, then Eqs. (5.24) and (5.25) apply.

Figure 5.3 shows the effect on material strength when $P > P_C$. The relaxation to Y_0 is achieved by assuming the material behaves like a Maxwell solid. The stress deviator S_{ij} is given by

$$s_{ij}^{n+1} = \hat{s}_{ij}^{n+1} \left[1 - \frac{\Delta t}{\delta} \left(1 - \frac{Y_{\varrho}}{J} \right) \right] \qquad J > Y_{\varrho}$$

$$s_{ij}^{n+1} = \hat{s}_{ij}^{n+1} \qquad J \leq Y_{\varrho}$$
(5.26)

where \hat{S}_{ij}^{n+1} and J are determined using Eqs. (5.10) and (5.12). The relaxation time, $\delta/(1-Y_j/J)$, increases as J approaches Y. This gives a smooth transition between the Maxwell solid and elastic behavior.

The behavior of $F(P_p)$ and the subsequent relaxation process given by Eqs. (5.22) through (5.26) is certainly an oversimplification of the role of pore fluid pressure in determining effective stress. Future work on this topic should merge experimental data of the type obtained by Duba, et al. [12] with the theoretical work of Garg and Nur. [11] We do feel, however, that the main features of effective stress are presently contained in the model. The primary input parameters which control the process are Y_0 , P_C and δ . At least some insight concerning the appropriate values for Y_0 and P_C should be able to be obtained from laboratory experiments. The value assumed for δ controls the relaxation time and we have used close—in time history data to determine this parameter.

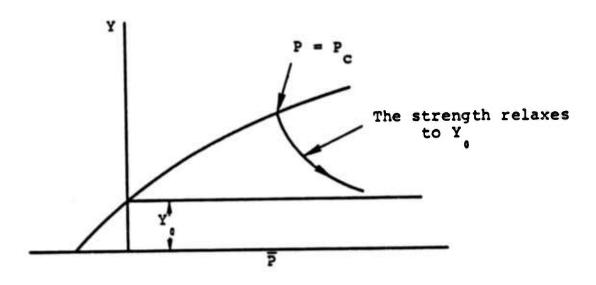


Figure 5.3. The relaxation of material strength when all air filled porosity has been removed.

5.6 SUMMARY OF INPUT PARAMETERS FOR MATERIAL STRENGTH

The limiting value of material strength, $Y(\overline{P})$, is obtained by specifying Y_0 , Y_m , \overline{P}_m and e_m . The function $Y(\overline{P})$ is given in Eq. (5.11). Except for the definition of \overline{P} , these parameters and $Y(\overline{P})$ are identical to that of Section 2.2. The overburden pressure P_0 is required to determine the ambient stress state of the material which is also required in Section 2.2. The crush pressure, P_C , plays a dual role. First it determines the limiting pressure at which the steepest release path occurs in the P-V plane, as shown in Fig. 2.4. Second, it establishes the pressure at which an effective stress law is used to determine material strength as shown in Fig. 5.3. The parameter δ determines the relaxation time of the stress deviators during the application of the effective stress law.

It is worth emphasizing that the only new parameter introduced in this section is δ , the relaxation time. However, the interpretation of material failure has changed considerably over the simple plastic flow model used in Section 2, Eq. (2.16). We now present some of the consequences of this revision.

5.7 THE CONSEQUENCES OF MATERIAL FAILURE ON THE EQUIVALENT ELASTIC SOURCE

A number of equivalent source calculations were performed using the material strength constitutive model presented in Sections 5.2 through 5.5. Table 5.1 lists the key material parameters used in the calculations along with the calculated values of $\Psi(\infty)$ and the final cavity radius, R_{cav} , for a device yield of 0.02 kt.

The input data listed in Table 5.1 has been taken from a number of sources. In particular the rock mechanics laboratory data of Stephens, Schock and Heard at the Lawrence Livermore Laboratory were used for granite and rhyolite. The data taken by Terra Tek for many events at NTS Area 12 were used for tuff. We should note, however, that no material strength data could be found for saturated rhyolite. We assumed that the material strength of this rock type would be the same as granite.

Figures 5.4 and 5.5 show velocity time histories recorded at two distances from Piledriver, a 60 kt event in granite at NTS (Perret [13]). In the past, most calculations using measured granite strength data have been unsuccessful in duplicating these velocity histories.

Figures 5.6 and 5.7 show the velocity histories obtained from our calculation of the Piledriver event (Calculation 130) and Table 5.2 summarizes the comparison between the calculation and the data. Cavity radius and peak velocities at the two stations agree with the data to within 5 percent. However, residual displacement at our first station is too high. This is due largely to the difficulty in reproducing the large negative phase of the velocity pulse of Fig. 5.4. Since this difficulty has been encountered in all previous calculations of Piledriver, it was decided to concentrate on the second station. Here we attained excellent agreement with the velocity data, including the long duration negative phase of the pulse. This gave us good agreement with the residual displacement measurement as well.

TABLE 5.1

REPRESENTATIVE MATERIAL PROPERTIES DATA FOR VARIOUS TEST AREAS

Name	gm/cm ^a	gm/cm³ km/sec	kbars 'p	kbars Y	kbars kbars kbars bars & kbars bars & m' m'' 'p' 'p' 'P' 'P' 'P' 'R'' 'R'' 'R''' 'R''' 'B'''' 'B''''' 'B''''''''	kbars	Para	•••	kbars Pc	Pare e	•4	* Î	2 0
Piledriver 130	2.67	5.33	207	7.7	0.3	9.0	0		0 1.0	•	•	0 0 9.23 3.2	3.2
Pahute Mesa Saturated Rhyolite 124 125	2.45	7	169	1.1	7.7 0.3 9.0 200	•	200	•	2.0	•	0 4.0 7.2	4.2	3.4
Area 12 Tuff 127	1.91	2.4	9		0.35 0.05 2.0	2.0		56 1.6	0.5	20	20 17.0 14.5	14.5	6.1
Hoggar Granite 129 128	2.67 Calcula	2.67 5.33 207 7.7 0.3 9.0 80 Calculation 128 used the same material properties as 129 except no tension failure was permitted in 128.	207 used the	7.7 scre m	0.3 aterial pr d in 128.	9.0 propertie	2:		pt 0	•	•	0 0 1.6	1.9

TABLE 5.2

COMPARISON BETWEEN DATA AND CALCULATION FOR PILEDRIVER

	Data	Calculation
Cavity radius (meters)	44.5	46.5
Peak Velocity ft/sec		
Station 1 (660' range)	110	103
Station 2 (1543 range)	18.8	19.4
Residual displacement (inches)		
Station 1	22	38
Station 2	7.2	6.1

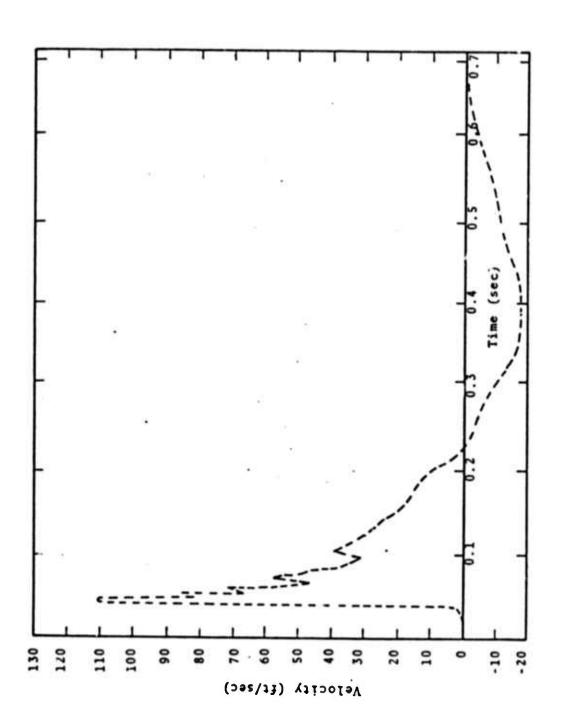
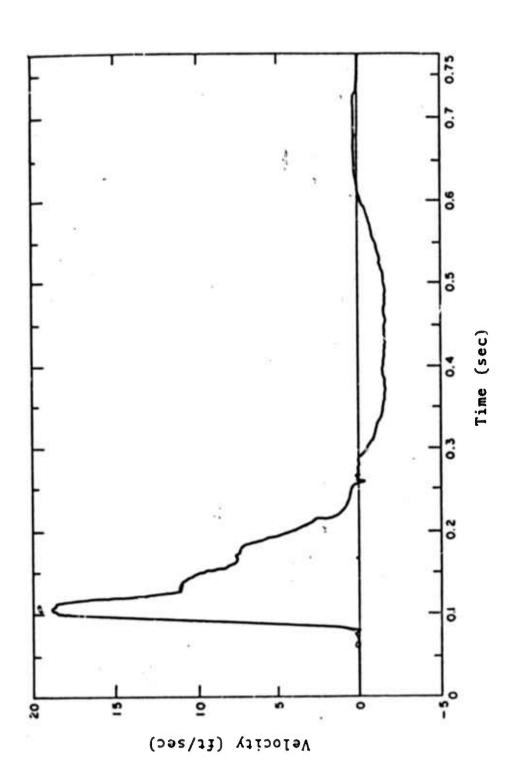


Figure 5.4. Velocity history at 668 feet horizontal range from Piledriver. [13]



Velocity history at 1543 feet horizontal range from Piledirver. [13] Figure 5.5.

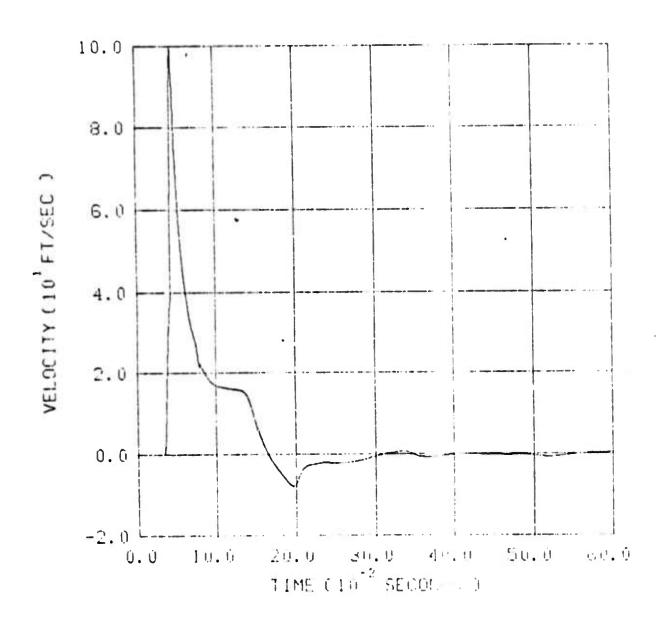


Figure 5.6. Calculated particle velocity at 668 feet from Piledriver.



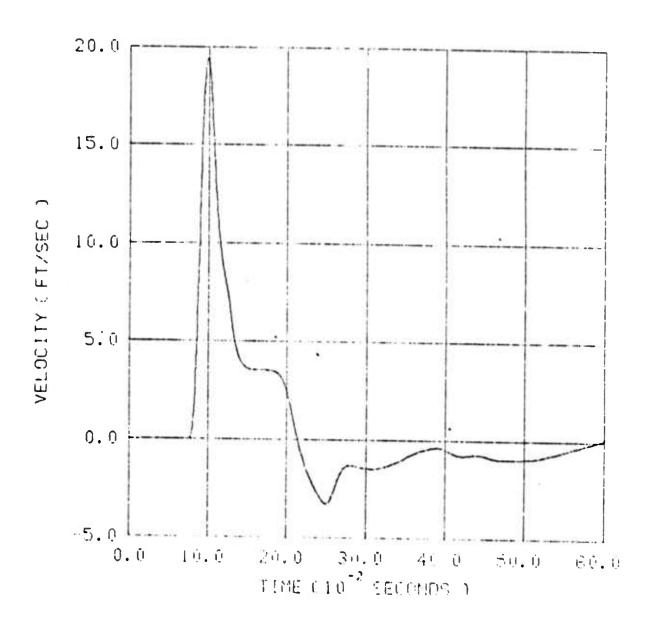


Figure 5.7. Calculated particle velocity at 1548 feet from Piledriver.

The Hoggar Granite calculation was intended to be representative of the source coupling for the French nuclear tests conducted in the Sahara. For this environment $P_{\rm c}$ was assumed infinite, a value which corresponds to a dry material. From Table 5.1 we find that the calculated cavity radius for Piledriver was 3.2 meters while for the Hoggar granite the cavity radius was 1.9 meters. This gives a factor of five difference in cavity volumes between the two test areas which is consistent with the published French test results.

Figure 5.8 shows the source functions corresponding to Calculations 127, 130 and 125. Both the Piledriver and Pahute Mesa spectra are noticeably peaked. The amplitude of the peak seems to correlate directly with Y.

In order to identify the physical phenomenon responsible for the peak we show the two Hoggar source spectra corresponding to Calculations 128 and 129 in Figs. 5.9 and 5.10. For Calculation 128 neither stress relaxation or tension failure was allowed and the source spectrum is flat; a result which is identical to the spectra shown in Fig. A.2 in Appendix A. Tension failure was permitted in Calculation 129 and the spectral peak for this calculation is approximately a factor of three greater than $\Psi(\infty)$.

The effect of stress relaxation on spectral shape is obtained by comparing the source spectra from Calculations 124 and 125. The total relaxation is less for 124 than for 125 since the corresponding values of P are 0.5 kbar and 1.0 kbar. Figure 5.11 compares the source spectra from these two calculations and indicates that the greater the stress relaxation the larger the spectral peak when compared to $\Psi(\infty)$.

Both tension failure and stress relaxation within the inelastic region seem to be the primary physical mechanisms responsible for producing a peaked spectrum. There is a good explanation for this since both mechanisms introduce a

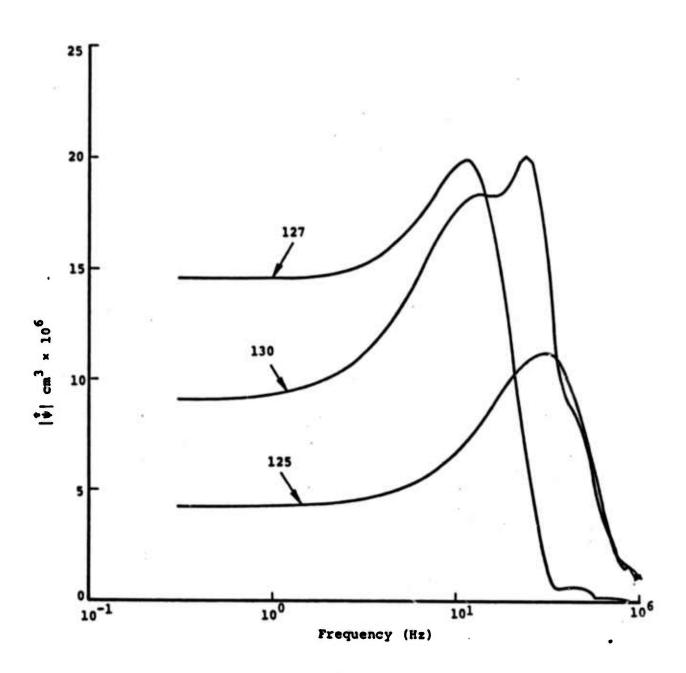
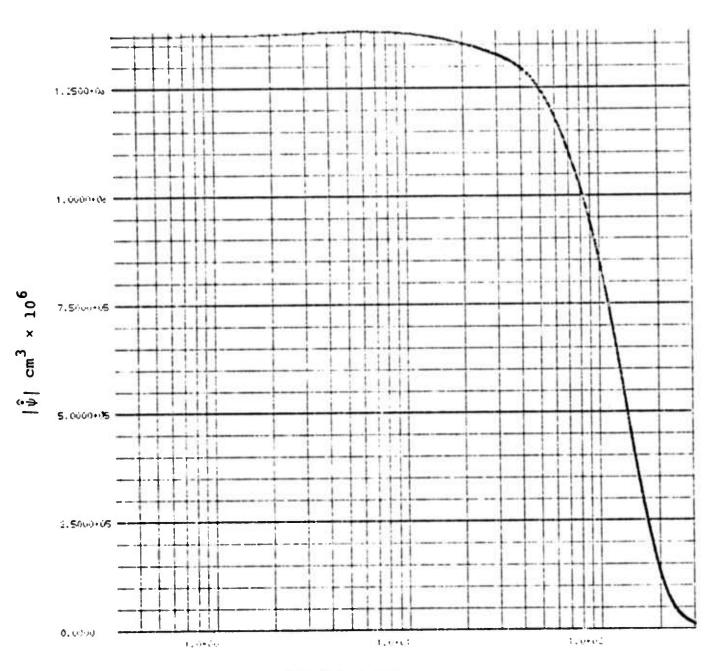
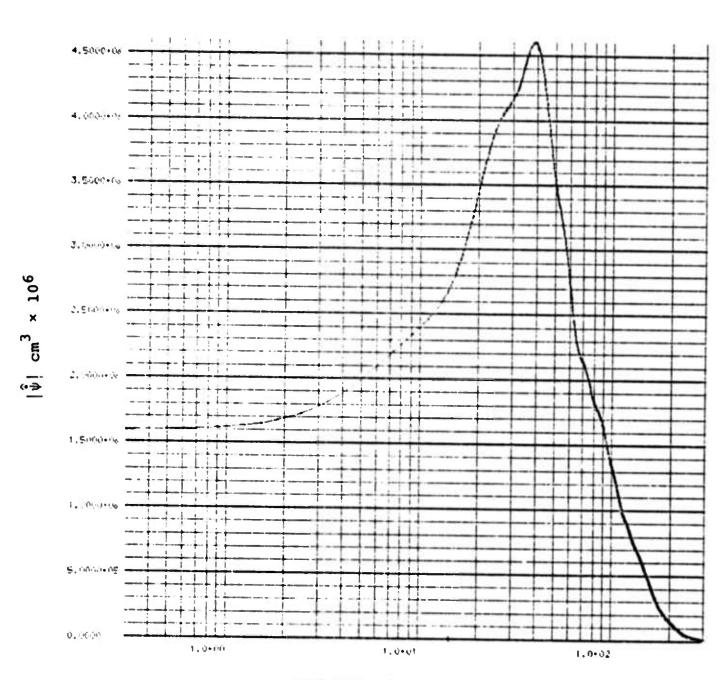


Figure 5.8. Equivalent source functions for Calculations 125, 127 and 130.



FREQUENCY (HZ)

Figure 5.9. Equivalent source function for Calculation 128.



FREQUENCY (HZ)

Figure 5.10. Equivalent source function for Calculation 129.

discontinuity in the magnitude of allowable stress between the nonlinear and linear elastic regions. This discontinuity causes the velocity field in the elastic region to rarefy, i.e., to become negative. The negative portion of the velocity time history, shown in Fig. 5.7, causes the peak in the source spectrum.

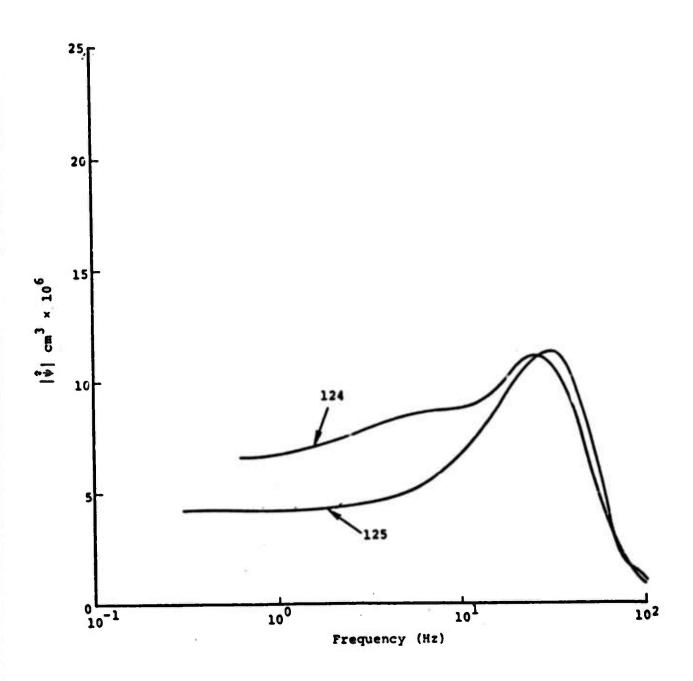


Figure 5.11. Equivalent source spectra for Calculations 124 and 125.

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APPENDIX A

THEORETICAL SCALING OF m_b WITH REDUCED DISPLACEMENT POTENTIAL

The equivalent elastic source for a spherically symmetric explosion is often represented by a reduced displacement potential, $\Psi(\tau)$. This quantity completely characterizes the coupling of the explosion into elastic waves in the immediate source region. Therefore, we expect proportionality between Ψ and teleseismic waves. While this proportionality can be quite complicated, it is useful to have a simple scaling law which indicates the variation of teleseismic body wave amplitude with Ψ and the properties of the emplacement material. Such a scaling relationship is derived in this appendix.

We begin by reviewing the most elementary model, that of a delta function source in a one material or homogeneous earth model. We then add material inhomogeneity in the form of a second material and allow a more complex source representation, arriving at the scaling

$$|u_t| \approx \alpha_s \Psi_\infty^e$$
 (A.1)

where $|u_t|$ is the amplitude of some consistently measured phase on the teleseismic record and Ψ_{∞}^{e} is a constant which characterizes the amplitude of the explosion generated pulse.

A Spherically Symmetric Source in a Homogeneous Earth

A spherically symmetric point source of dilatational waves can be characterized by a reduced displacement potential. That is, the radial displacement $u_{\bf g}({\bf R},t)$ can be written

$$u_s(R,t) = -\frac{\partial}{\partial R} \left(\frac{\Psi(\tau)}{R} \right) = \frac{\Psi(\tau)}{R^2} + \frac{\dot{\Psi}(\tau)}{R\alpha_s}$$
 (A.2)

where $\tau = t - R/\alpha_s$, the time retarded by the P wave travel time to the position R. Here α_s is the P wave velocity in the source region.

Taking Fourier transforms, (A.2) may be written

$$\hat{\mathbf{u}}_{\mathbf{S}}(\mathbf{R},\omega) = \left(\frac{1}{\mathbf{R}^2} + \frac{i\mathbf{k}_{\alpha}}{\mathbf{R}}\right) \hat{\boldsymbol{\Psi}}(\omega) e^{-i\mathbf{k}_{\alpha}\mathbf{R}}, \qquad (A.3)$$

where $k_{\alpha} \equiv \omega/\alpha_s$. If $k_{\alpha}R >> 1$, that is, if R is many wavelengths, only the "far field" terms need to be retained and

$$\hat{\mathbf{u}}_{\mathbf{g}}(\mathbf{R},\omega) = \frac{\hat{\mathbf{v}}(\omega)}{\mathbf{R}\mathbf{u}_{\mathbf{g}}} e^{-\mathbf{i}\mathbf{k}_{\alpha}\mathbf{R}} . \tag{A.4}$$

For teleseismic body wave amplitudes only frequencies less than about 2.5 Hz are of importance. Then for low yield explosions $\hat{\Psi}(\omega)$ is essentially constant and equal to Ψ_{∞} , the static value of the reduced displacement potential. Therefore, for a low yield source of the type discussed in Section IV, and a homogeneous earth model, we have

$$|u_t| \approx \frac{\Psi_{\infty}}{\alpha_g}$$
 (A.5)

For higher yields the teleseismic frequency band includes portions of the peak on the $\hat{\Psi}(\omega)$ curves. Therefore Eq. (A.5) is improved by replacing Ψ_{∞} by an "equivalent" value, Ψ_{∞}^{e} , which is yield dependent. Once again, this is

The assumption that $\Psi(\omega) \approx \Psi_{\infty}$ for the teleseismic frequencies is equivalent to assuming that the source appears to be a delta function of amplitude Ψ_{∞} to a teleseismic observer.

equivalent to assuming that the source behaves, teleseismically, as if it were a delta function of amplitude Ψ_{∞}^{e} . Computation of theoretical seismograms show that this assumption is not bad. The value of Ψ^{e} is taken to be the value of $\Psi(\omega)$ at the frequency which cube root scales to the yield of interest. With this modification, (A.5) is the appropriate scaling law for the sources of Section IV in a homogeneous earth model.

A Spherically Symmetric Source in a Two Layered Whole Space

Let us now assume that the spherically symmetric source is located in a space composed of two layers. The geometry is shown in Fig. A.1.

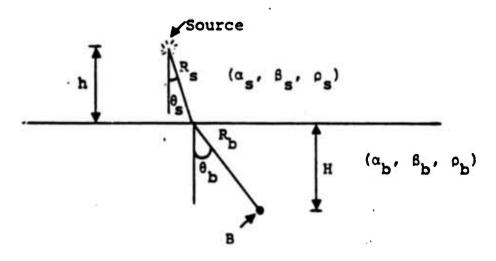


Figure A.1. A source in a two layered medium.

If attention is restricted to the far field, the displacement in potential form is given by (A.4). Then the P wave displacement at the point B below the interface is given by

$$\hat{\mathbf{u}}_{\mathbf{b}} = \frac{\mathbf{A}}{\hat{\mathbf{L}}} \frac{\hat{\Psi}(\mathbf{\omega})}{\alpha_{\mathbf{c}}} e^{-\mathbf{i}\boldsymbol{\omega}\tau_{\mathbf{b}}} , \qquad (A.6)$$

where

$$\begin{split} & L^{2} = \left(R_{s} + \frac{\alpha_{b}}{\alpha_{s}} R_{b}\right) \left(R_{s} + \frac{\cos^{2}\theta_{s}}{\cos^{2}\theta_{b}} \frac{\alpha_{b}}{\alpha_{s}} R_{b}\right), \\ & A = \frac{2\rho_{s}\alpha_{s}}{D} P_{1} \left(\beta_{b}P_{2}X + \beta_{s}P_{s}Y\right), \\ & D = \alpha_{s}\alpha_{b}\beta_{s}\beta_{b} \frac{Z^{2}}{c^{2}} + \alpha_{b}\beta_{b}P_{1}P_{2}X^{2} + \alpha_{s}\beta_{s}P_{1}P_{s}Y^{2}, \\ & + \rho_{s}\rho_{b} \left(\beta_{s}\alpha_{b}P_{1}P_{s} + \alpha_{s}\beta_{b}P_{2}P_{3}\right) + \frac{q^{2}}{c^{2}} P_{1}P_{2}P_{s}P_{s}, \\ & X = \rho_{b} - \frac{q}{c^{2}}, Y = \rho_{s} + \frac{q}{c^{2}}, Z = \rho_{b} - \rho_{s} - \frac{q}{c^{2}}, \\ & P_{1} = \cos\theta_{s}, P_{2} = \left(1 - \frac{\beta_{s}^{2}}{c^{2}}\right)^{1/2}, P_{3} = \cos\theta_{b}, P_{4} = \left(1 - \frac{\beta_{b}^{2}}{c^{2}}\right)^{1/2}, \\ & C = \frac{\alpha_{s}}{\sin\theta_{s}} = \frac{\alpha_{b}}{\sin\theta_{b}}, q = 2 \left(\rho_{b}\beta_{b}^{2} - \rho_{s}\beta_{s}^{2}\right). \end{split}$$

In (A.6) the exponential term is merely the phase shift due to the travel time from the source to B. The L is the spreading term for the spherical wave. For a homogeneous material, L = R, the distance between the source and B. The coefficient A is the refraction coefficient which depends on the properties of the two media and the angle of incidence which we denote θ_s .

Application to Teleseismic Body Waves

The two layered model of Fig. A.l is more realistic than a homogeneous earth model for illuminating the gross

properties of the teleseismic body waves excited by explosions in different source materials. For example, consider the NTS explosions. The resulting teleseismic body waves observed at a single seismograph station are expected to vary due to differences only in the explosion source and near source crustal structure. All other features of the travel path are essentially held constant.

Below some depth one expects all of the NTS to be underlain by a common material. The material denoted by a subscript b can be thought of as this basement rock. Explosions at NTS have been detonated in such widely varying materials as alluvium, tuff, rhyolite and granite. These are characterized by P wave velocities ranging from 2.0 - 5.5 km/sec. The subscript s can be taken to represent this source material. For the long wavelengths (> 2 km) of interest for teleseismic waves, the single interface transition between the source and basement materials is a useful approximation.

The body waves reaching the teleseismic field exit the source region at near vertical takeoff angles. That is, $\cos\theta_{\rm S} \approx 1$ and $\cos\theta_{\rm b} \approx 1$. Then it is clear from (A.7) that

$$L \approx R_s + \frac{\alpha_b}{\alpha_s} R_b . \tag{A.8}$$

However, since the basement rock is held in common from event to event, one can choose R_b arbitrarily large while R_s remains rather small. Thus, comparing different events at NTS, $L \approx \alpha_b/\alpha_s$.

The complexity of its form makes it difficult to deduce a similarly simple relationship for A. However, numerical experiments using values for typical geologic materials and horizontal wave speed c (or takeoff angle θ_b) show that A is very nearly proportional to α_s/α_b . Taking c = 12 km/sec and α_b = 6.0 km/sec, β_b = 3.5 km/sec, ρ_b = 2.8 gm/cm³, a number of illustrative examples are given in Table A.1.

TABLE A.1. THE REFRACTION COEFFICIENT A FOR TYPICAL SOURCE MATERIALS

α _s	β _s	Ps	θ_s	A	$\frac{\alpha_{b}^{A/\alpha}s}$
2.40	1.30	1.86	11.53	0.4199	1.05
2.80	1.40	1.95	13.29	0.4906	1.05
3.30	1.90	2.25	15.66	0.6130	1.11
4.00	2.50	2.40	19.26	0.7273	1.09
4.50	2.65	2.60	21.71	0.8211	1.09
5.10	2.75	2.70	24.20	0.9009	1.06
5.50	3.10	2.75	26.74	0.9475	1.03

For short period teleseismic body waves $\Psi(\omega)$ may be replaced by the constant $\Psi_{\infty}^{\mathbf{e}}$. Then combining all these approximations with (A.6), we find that teleseismic amplitude is proportional to $\alpha_{\mathbf{g}}\Psi_{\infty}^{\mathbf{e}}$. That is, for the two layered earth model we deduce that

$$|u_t| \approx \alpha_s \Psi_\infty^e$$
 (A.9)

It should be emphasized that it is the direct P wave that scales with $\alpha_g \Psi_\infty^e$. Extending this to scale m_b implicitly assumes that the influence of the free surface reflected phase and details of the near source geology are invarient between events. The usefulness of (A.9) is degraded to the extent that this is not true.

Calculation of the RDP

The RDP is calculated in the following manner. If we assume that the displacement is linear within a given time interval defined by

$$\tau_1 < \tau < \tau_2$$

then Eq. (A.2) gives

$$\Psi(\tau_2) = aR^2 \left[\tau_2 - \tau_1 e^X\right] + R^2 \left(b - \frac{aR}{ag}\right) \left[1 - e^X\right] + e^X \Psi(\tau_1) \quad (A.10)$$

where

$$a = \frac{u(\tau_1, R) - u(\tau_2, R)}{\tau_1 - \tau_2}$$

$$b = \frac{-u(\tau_1, R)\tau_2 + u(\tau_2, R)\tau_1}{\tau_1 - \tau_2}$$

$$x = \frac{\alpha_{s}(\tau_{1} - \tau_{2})}{R}$$

At late times when the displacement becomes constant, $\tau_{_{2}}$ can be much larger than $\tau_{_{1}}$ so that

$$b = u(\infty, R)$$

The steady-state value of the RDP becomes

$$\Psi(\infty) = R^2 \ u(\infty, R) \tag{A.11}$$

This result also follows directly from Eq. (A.2) since

$$\dot{\Psi}(\infty) = 0 \tag{A.12}$$

In an entirely analogous manner, the reduced velocity potential (RVP) may be computed from the velocity-time history of a particle. If $\dot{u}(t,R)$ is radial particle velocity, then

$$\dot{\mathbf{u}}(\mathsf{t},\mathsf{R}) = -\frac{\partial}{\partial \mathsf{R}} \frac{\dot{\mathbf{v}}(\mathsf{\tau})}{\mathsf{R}} = \frac{\dot{\mathbf{v}}(\mathsf{\tau})}{\mathsf{R}^2} + \frac{\ddot{\mathbf{v}}(\mathsf{\tau})}{\alpha_{\mathsf{s}}\mathsf{R}} \tag{A.13}$$

where $\Psi(\tau)$ is the RVP.

If $\dot{\Psi}(\tau)$ replaces $\Psi(\tau)$ and $\dot{u}(\tau,R)$ replaces $u(\tau,R)$ in Eq. (A.10), then the algorithm for calculating the RVP follows directly.

It is highly desirable to transform the equivalent source into the frequency domain in order to identify the frequency dependence of the spectrum required for equivalent source comparison. Since $\mathring{\Psi}(\infty)=0$, then $\mathring{\Psi}$ is bounded at zero frequency and the Fourier transform of $\mathring{\Psi}(\tau)$ may be done numerically.

An important relation is that the zero frequency limit of the RVP is equal to the steady state value of the RDP. Since

$$\hat{\Psi}(f=0) = \int_{0}^{\infty} \Psi(\tau) d\tau = \int_{0}^{\infty} \frac{d\Psi(\tau)}{d\tau} d\tau$$

Therefore, since $\Psi(0) = 0$,

$$\hat{\Psi}(U) = \Psi(\infty) \tag{A.14}$$

Reduced velocity potentials have been computed for a variety of near source geologic environments. These calculations were performed on a Lagrangian finite difference code,

SKIPPER (see Appendix B), that simulates a propagating stress wave in one space dimension. The difference equations in the code are identical to those given by Cherry and Petersen. [7] The code is capable of carrying the propagating stress wave into the small displacement, elastic regime and yet flexible enough to permit appropriate material response formulations in the large displacement, nonlinear regime.

Figure A.2 shows $|\Psi|$ versus frequency for a nuclear explosion of 20 kt. Each source function in the figure is identified by an integer and corresponds to a 20 kt explosion detonated in a geologic environment with a specific set of material properties given in Section III. For all practical purposes the spectra are flat within the teleseismic frequency band (f < 3 Hz) for yields up to at least 200 kt. This implies that teleseismic body wave and surface wave amplitudes should be directly proportional to the yield of the explosion. Deviation from this linear scaling should occur within the yield interval 200 kt < W < 1,000 kt, depending on the source function used to represent the rock material near the source.

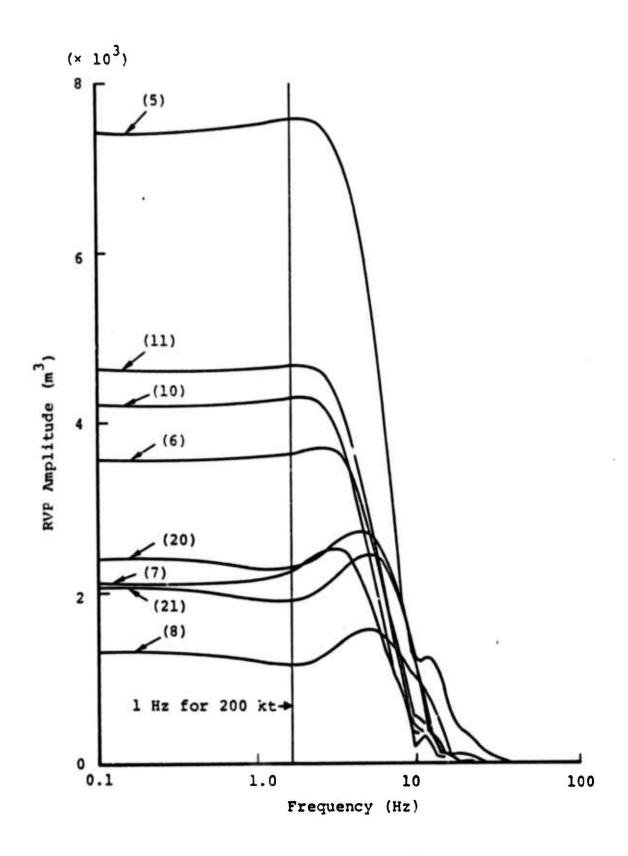


Figure A.2. Source functions scaled to 20 kt.

APPENDIX B

THE SKIPPER CODE

B.1 DESCRIPTION OF THE CODE

SKIPPER is a Lagrangian, finite difference code that simulates a propagating stress wave in one space dimension. The code is capable of carrying the propagating stress wave into the small displacement, elastic regime and yet flexible enough to permit appropriate material response formulations in the large displacement, nonlinear regime.

A schematic of one computational cycle is shown in Fig. B.l. The equation of motion for spherical symmetry is

$$\rho \dot{\mathbf{u}} = -\left[\frac{\partial \sigma}{\partial \mathbf{r}} + \frac{2(\sigma - \sigma)}{\mathbf{r}}\right] \tag{B.1}$$

where ρ is density, u is particle velocity, u is particle acceleration and σ and σ are the radial and tangential components of the stress tensor with compression positive.

Strain rates in the radial and tangential direction are

$$\dot{\varepsilon}_{1} = \frac{\partial u}{\partial r} \tag{B.2}$$

$$\frac{\cdot}{\varepsilon_2} = \frac{u}{r}$$
 (B.3)

The stress tensor may be decomposed into an isotropic component (P) and a deviatoric component (S $_1$ and S $_2$) where

$$\sigma_{i} = P - S_{i} \tag{B.4}$$

Due to symmetry

$$S_1 = -2S_2 \tag{B.5}$$

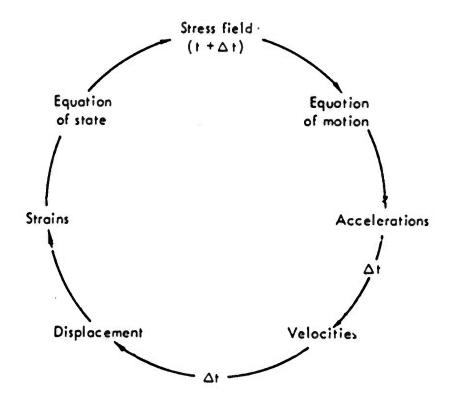


Figure B.1. Cycle of interactions treated in calculating stress wave propagation. Taken from Ref. 5.

Hooke's law gives

$$\dot{P} = -\kappa \frac{\dot{v}}{v} \tag{B.5}$$

$$\dot{S}_{1} = \frac{4}{3} \mu (\dot{\varepsilon}_{1} - \dot{\varepsilon}_{2}) \tag{B.7}$$

where k and μ are the bulk and shear moduli, the dot indicates a time derivative along a particle path and from conservation of mass the rate of change of volume is

$$\frac{\dot{\mathbf{v}}}{\mathbf{v}} = \dot{\boldsymbol{\varepsilon}}_1 + 2\dot{\boldsymbol{\varepsilon}}_2 \tag{B.8}$$

For large and/or energy dependent deformations Eq. (B.6) is replaced by

$$P = f_1(v,e) \text{ or } P = g(\overline{n},e)$$
 (B.9)

where v is specific volume.

The specific internal energy (e) is obtained from

$$\rho \dot{\mathbf{e}} = -P \dot{\mathbf{v}} + S_1 \dot{\varepsilon}_1 + 2S_2 \dot{\varepsilon}_2 \tag{B.10}$$

and the total volumetric strain $(\overline{\eta})$ is defined to be

$$\overline{\eta} = \frac{v^0 - v}{v^0} = 1 - \frac{v}{v_0} = 1 - \frac{\rho_0}{\rho}$$
 (B.11)

The difference equation used in the code that corresponds to Eqs. (B.1), (B.2), (B.3), (B.10) and (B.11) are the same as those used by Cherry and Petersen. [7]

Basically, Eqs. (B.7) and (B.9) should be regarded as an initial attempt by the code to determine P and S at the current time cycle. Modification of these stresses then depend on the constitutive relation used in the equation of state section of the code.

An artificial viscosity is used in the momentum and energy equations [(B.1) and (B.10)) and has the same form as that reported by Wilkins. [8] This quantity replaces discontinuous jumps in particle velocity in the calculation with a rapid but continuous change.

B.2 SKIPPER COMPARISON TO AN ANALYTIC SOLUTION

Normally, the only way to evaluate a numerical procedure is to examine its ability to produce the answer to a previously solved problem. Also, in order to calculate the equivalent elastic source function from a contained nuclear explosion, the calculations must be continued well into the small displacement, elastic regime. In order for the calculation to be meaningful in this regime the numerics in the code must not introduce errors in the solution when displacements are small.

A well-known analytic solution, commonly called Blake's solution, [9] to an elastic wave propagation problem has been used for our analytic comparison. The following parameters were used in the calculation:

Cavity Radius "10 meters

Density = 2 g/cc

Compressional Velocity = 5 m/msec

Bulk Modulus = 333.33 kbar

Shea: Modulus = 125 kbar

Cavity Load (kbar) = e^{-t} (t in msec)

Zone Size = 10 cm

Figures B.2 and B.3 show particle velocity versus radius at 1.32 msec and 6.81 msec. Figures B.4 and B.5 show displacement and mean stress versus time 10 m from the cavity wall.

The numerical results from SKIPPER agree quite well with the analytic solution. Displacement time histories of the type shown in Fig. B.4 provide the input for the RDP calculation.

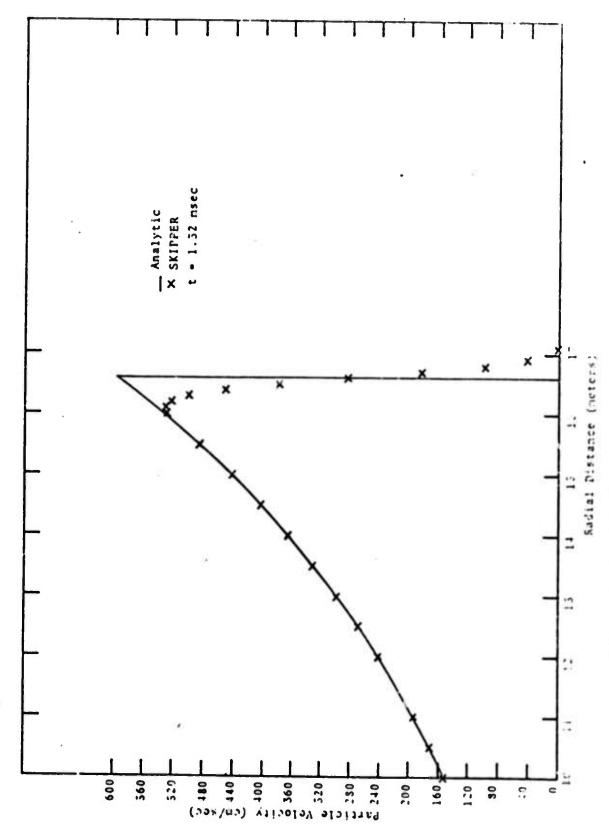
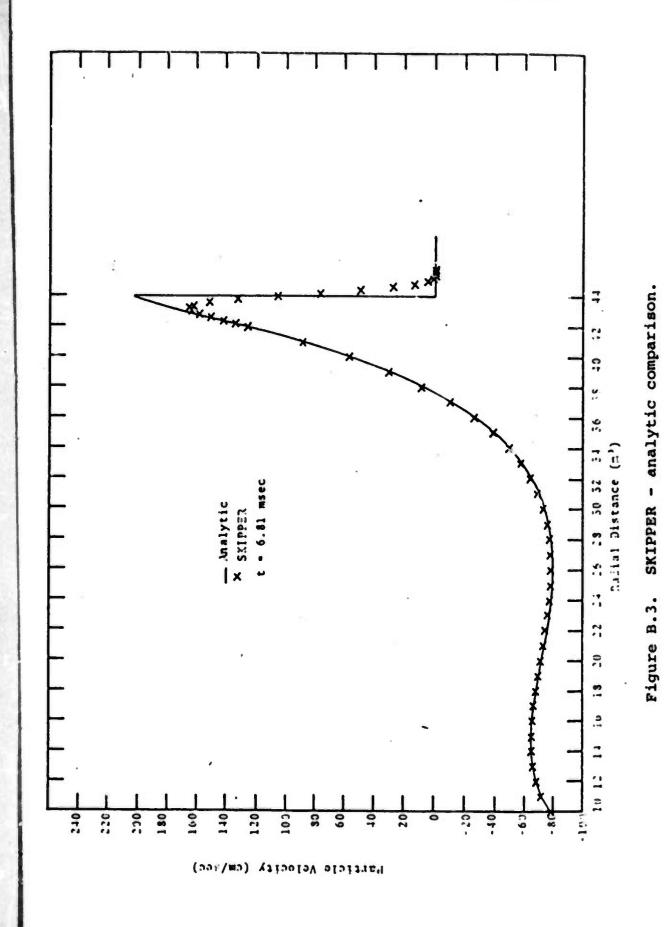
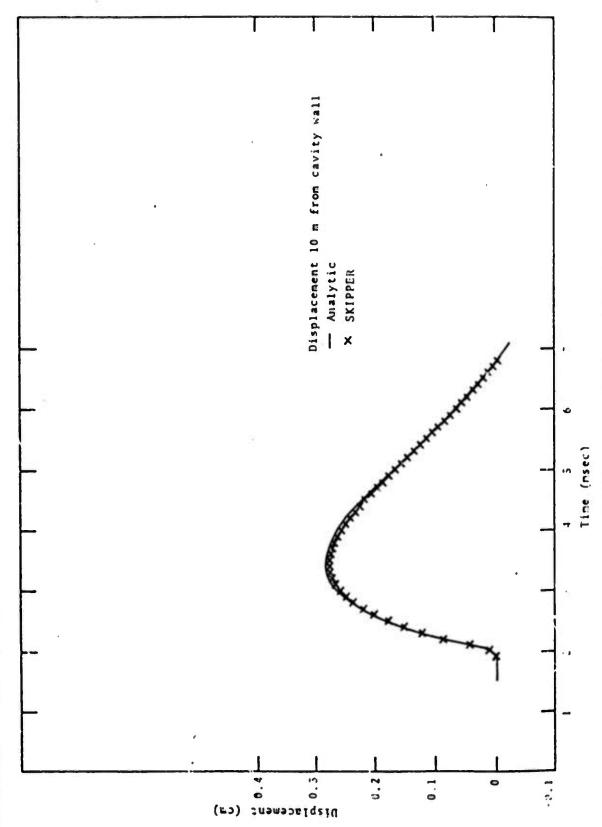


Figure B.2. SKIPPER - analytic comparison.



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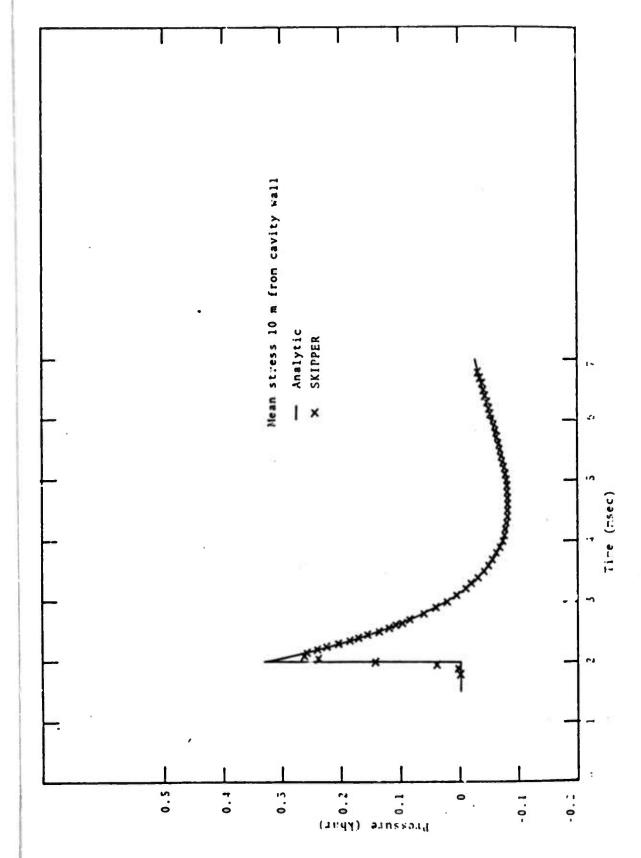


Figure B.5. SKIPPER - analytic comparison.